Full wwPDB X-ray Structure Validation Report (i)

Aug 16, 2023 – 01:21 PM EDT

PDB ID	:	2A2J
Title	:	Crystal structure of a putative pyridoxine 5'-phosphate oxidase (Rv2607) from Mycobacterium tuberculosis
Authors	:	Pedelacq, J.-D.; Rho, B.-S.; Kim, C.-Y.; Waldo, G.S.; Lekin, T.P.; Segelke, B.W.; Rupp, B.; Hung, L.-W.; Kim, S.-I.; Terwilliger, T.C.; Mycobacterium Tuberculosis Structural Proteomics Project (XMTB)
Deposited on	:	2005-06-22
Resolution	:	2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) were used in the production of this report:

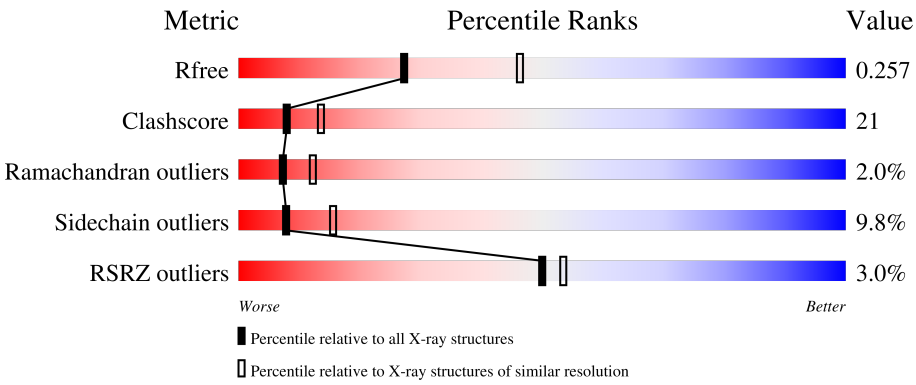
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	246	3% 58% 17% 6% • 17%
1	B	246	2% 55% 21% 6% • 17%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3281 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Pyridoxamine 5'-phosphate oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	203	1603	1007	291	299	6	0	0	0
1	B	203	1603	1007	291	299	6	0	0	0

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	cloning artifact	UNP P65682
A	-18	GLY	-	cloning artifact	UNP P65682
A	-17	SER	-	cloning artifact	UNP P65682
A	-16	SER	-	cloning artifact	UNP P65682
A	-15	HIS	-	expression tag	UNP P65682
A	-14	HIS	-	expression tag	UNP P65682
A	-13	HIS	-	expression tag	UNP P65682
A	-12	HIS	-	expression tag	UNP P65682
A	-11	HIS	-	expression tag	UNP P65682
A	-10	HIS	-	expression tag	UNP P65682
A	-9	SER	-	cloning artifact	UNP P65682
A	-8	SER	-	cloning artifact	UNP P65682
A	-7	GLY	-	cloning artifact	UNP P65682
A	-6	LEU	-	cloning artifact	UNP P65682
A	-5	VAL	-	cloning artifact	UNP P65682
A	-4	PRO	-	cloning artifact	UNP P65682
A	-3	ARG	-	cloning artifact	UNP P65682
A	-2	GLY	-	cloning artifact	UNP P65682
A	-1	SER	-	cloning artifact	UNP P65682
A	0	HIS	-	cloning artifact	UNP P65682
A	29	CME	CYS	modified residue	UNP P65682
A	225	GLY	-	cloning artifact	UNP P65682
A	226	SER	-	cloning artifact	UNP P65682
B	-19	MET	-	cloning artifact	UNP P65682
B	-18	GLY	-	cloning artifact	UNP P65682

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	SER	-	cloning artifact	UNP P65682
B	-16	SER	-	cloning artifact	UNP P65682
B	-15	HIS	-	expression tag	UNP P65682
B	-14	HIS	-	expression tag	UNP P65682
B	-13	HIS	-	expression tag	UNP P65682
B	-12	HIS	-	expression tag	UNP P65682
B	-11	HIS	-	expression tag	UNP P65682
B	-10	HIS	-	expression tag	UNP P65682
B	-9	SER	-	cloning artifact	UNP P65682
B	-8	SER	-	cloning artifact	UNP P65682
B	-7	GLY	-	cloning artifact	UNP P65682
B	-6	LEU	-	cloning artifact	UNP P65682
B	-5	VAL	-	cloning artifact	UNP P65682
B	-4	PRO	-	cloning artifact	UNP P65682
B	-3	ARG	-	cloning artifact	UNP P65682
B	-2	GLY	-	cloning artifact	UNP P65682
B	-1	SER	-	cloning artifact	UNP P65682
B	0	HIS	-	cloning artifact	UNP P65682
B	29	CME	CYS	modified residue	UNP P65682
B	225	GLY	-	cloning artifact	UNP P65682
B	226	SER	-	cloning artifact	UNP P65682

- Molecule 2 is water.

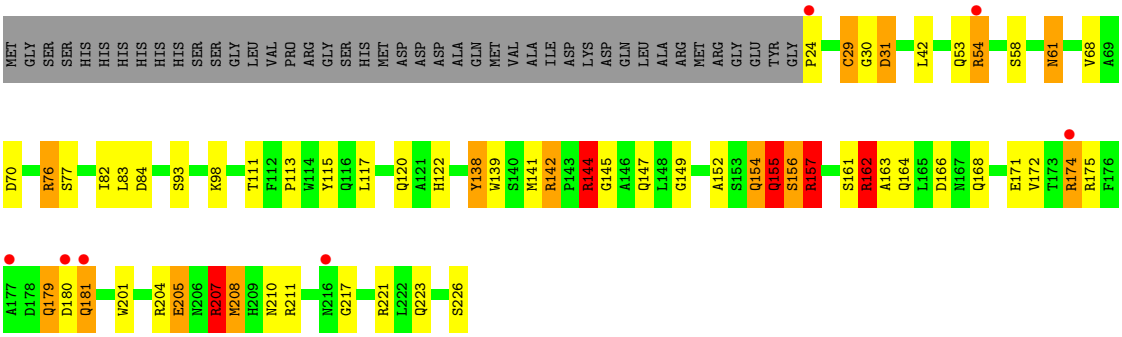
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	44	44	44	0	0
			Total	O		
2	B	31	31	31	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

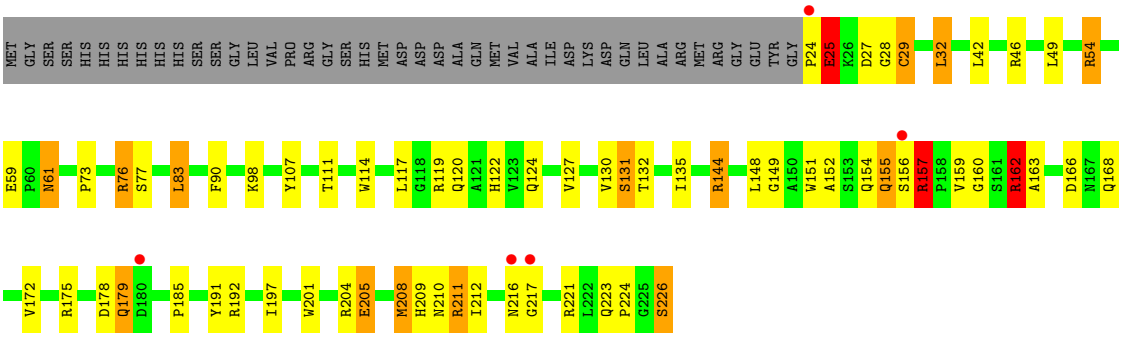
• Molecule 1: Pyridoxamine 5'-phosphate oxidase

Chain A:



• Molecule 1: Pyridoxamine 5'-phosphate oxidase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	69.61Å 91.11Å 92.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.48 – 2.50 29.48 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.48-2.50) 100.0 (29.48-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.75 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.209 , 0.260 0.209 , 0.257	Depositor DCC
R_{free} test set	1535 reflections (7.29%)	wwPDB-VP
Wilson B-factor (Å^2)	37.3	Xtriage
Anisotropy	0.607	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 30.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.020 for -h,l,k	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	3281	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.26% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CME

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\|>5$	RMSZ	$\#\|Z\|>5$
1	A	1.13	6/1634 (0.4%)	1.13	16/2220 (0.7%)
1	B	1.14	6/1634 (0.4%)	1.12	10/2220 (0.5%)
All	All	1.14	12/3268 (0.4%)	1.13	26/4440 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
All	All	0	3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	107	TYR	CD2-CE2	9.66	1.53	1.39
1	A	205	GLU	C-N	7.41	1.51	1.34
1	A	138	TYR	CG-CD1	7.31	1.48	1.39
1	B	107	TYR	CD1-CE1	7.28	1.50	1.39
1	A	30	GLY	C-N	-6.47	1.19	1.34
1	B	205	GLU	CB-CG	6.47	1.64	1.52
1	A	174	ARG	C-N	-6.42	1.19	1.34
1	A	138	TYR	CE2-CZ	5.85	1.46	1.38
1	A	161	SER	C-N	-5.63	1.21	1.34
1	B	160	GLY	CA-C	5.37	1.60	1.51
1	B	178	ASP	CG-OD1	5.19	1.37	1.25
1	B	131	SER	C-N	-5.04	1.22	1.34

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	54	ARG	NE-CZ-NH2	-14.41	113.09	120.30
1	B	54	ARG	NE-CZ-NH1	9.01	124.80	120.30
1	A	31	ASP	O-C-N	8.03	135.54	122.70
1	A	30	GLY	O-C-N	-7.96	109.97	122.70
1	A	157	ARG	NE-CZ-NH2	7.80	124.20	120.30
1	B	144	ARG	NE-CZ-NH2	7.50	124.05	120.30
1	A	175	ARG	NE-CZ-NH2	7.38	123.99	120.30
1	B	154	GLN	O-C-N	7.32	134.41	122.70
1	A	207	ARG	NE-CZ-NH2	7.30	123.95	120.30
1	B	157	ARG	NE-CZ-NH2	7.26	123.93	120.30
1	A	142	ARG	NE-CZ-NH2	7.12	123.86	120.30
1	B	211	ARG	NE-CZ-NH2	7.07	123.83	120.30
1	A	30	GLY	C-N-CA	6.91	138.98	121.70
1	A	144	ARG	O-C-N	-6.87	111.52	123.20
1	A	144	ARG	NE-CZ-NH2	6.42	123.51	120.30
1	B	208	MET	CG-SD-CE	6.25	110.19	100.20
1	A	208	MET	CG-SD-CE	6.19	110.10	100.20
1	A	141	MET	CG-SD-CE	6.11	109.98	100.20
1	A	31	ASP	CA-C-N	-5.69	104.67	117.20
1	B	162	ARG	NE-CZ-NH2	5.65	123.12	120.30
1	A	205	GLU	C-N-CA	-5.51	107.92	121.70
1	B	154	GLN	CA-C-N	-5.44	105.24	117.20
1	A	84	ASP	CB-CG-OD2	-5.33	113.51	118.30
1	A	162	ARG	NE-CZ-NH2	5.24	122.92	120.30
1	A	30	GLY	CA-C-N	5.12	128.47	117.20
1	B	119	ARG	NE-CZ-NH2	-5.09	117.76	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	29	CME	Mainchain
1	B	25	GLU	Peptide
1	B	28	GLY	Mainchain

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1603	0	1541	90	1
1	B	1603	0	1544	74	1
2	A	44	0	0	11	0
2	B	31	0	0	4	0
All	All	3281	0	3085	135	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 21.

All (135) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:208:MET:CE	1:B:152:ALA:CB	1.82	1.55
1:A:208:MET:HE1	1:B:152:ALA:CB	1.41	1.37
1:A:208:MET:CE	1:B:152:ALA:HB1	1.45	1.36
1:A:208:MET:CE	1:B:152:ALA:HB3	1.56	1.20
1:A:155:GLN:HE21	1:A:155:GLN:CA	1.57	1.14
1:A:155:GLN:HA	1:A:155:GLN:NE2	1.59	1.11
1:A:208:MET:HE3	1:B:152:ALA:HB1	1.08	1.04
1:A:154:GLN:NE2	1:A:157:ARG:HE	1.56	1.02
1:A:61:ASN:H	1:A:61:ASN:HD22	1.07	0.95
1:A:211:ARG:HH12	1:A:226:SER:HA	1.30	0.94
1:A:24:PRO:HD3	2:A:229:HOH:O	1.67	0.93
1:A:155:GLN:CA	1:A:155:GLN:NE2	2.21	0.93
1:A:162:ARG:HH11	1:B:210:ASN:HD21	1.05	0.93
1:A:208:MET:HE2	1:B:152:ALA:CB	1.97	0.91
1:A:155:GLN:HE21	1:A:155:GLN:HA	0.76	0.91
1:A:154:GLN:HE21	1:A:157:ARG:HE	1.11	0.91
1:A:162:ARG:NH1	1:B:210:ASN:HD21	1.71	0.89
1:A:210:ASN:HD22	1:A:221:ARG:HH12	1.16	0.88
1:A:211:ARG:NH1	1:A:226:SER:HA	1.87	0.88
1:B:211:ARG:NH1	1:B:226:SER:HA	1.89	0.88
1:A:154:GLN:HE21	1:A:157:ARG:NE	1.76	0.83
1:A:154:GLN:NE2	1:A:157:ARG:NE	2.26	0.82
1:B:155:GLN:HG3	1:B:156:SER:N	1.94	0.81
1:A:208:MET:HE1	1:B:152:ALA:HB3	0.81	0.80
1:B:211:ARG:HH12	1:B:226:SER:HA	1.47	0.76
1:A:61:ASN:HD22	1:A:61:ASN:N	1.83	0.73
1:A:154:GLN:O	1:A:155:GLN:O	2.08	0.72
1:A:210:ASN:HD22	1:A:221:ARG:NH1	1.89	0.70
1:A:120:GLN:HE22	1:B:77:SER:H	1.40	0.70
1:B:24:PRO:HD3	1:B:205:GLU:HB2	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:GLN:O	1:A:155:GLN:C	2.31	0.68
1:B:204:ARG:H	1:B:209:HIS:HD2	1.41	0.68
1:B:61:ASN:HD22	1:B:61:ASN:H	1.41	0.67
1:B:122:HIS:HE1	2:B:233:HOH:O	1.78	0.66
1:A:122:HIS:HE1	2:A:228:HOH:O	1.78	0.66
1:A:179:GLN:HE21	1:A:180:ASP:N	1.93	0.66
1:A:154:GLN:H	1:B:223:GLN:NE2	1.94	0.66
1:B:155:GLN:CG	1:B:156:SER:N	2.59	0.66
1:A:70:ASP:CB	2:A:266:HOH:O	2.44	0.64
1:A:179:GLN:NE2	1:A:181:GLN:H	1.97	0.64
1:A:70:ASP:HB3	2:A:266:HOH:O	1.97	0.63
1:A:168:GLN:O	1:A:172:VAL:HG23	1.98	0.62
1:A:111:THR:OG1	1:A:122:HIS:CD2	2.53	0.62
1:A:115:TYR:HB3	1:B:59:GLU:HG3	1.81	0.62
1:A:223:GLN:HE22	1:B:168:GLN:HE22	1.48	0.61
1:A:24:PRO:HD3	1:A:205:GLU:H	1.66	0.61
1:A:179:GLN:HE21	1:A:179:GLN:CA	2.14	0.61
1:A:179:GLN:HE21	1:A:180:ASP:H	1.47	0.61
1:A:61:ASN:H	1:A:61:ASN:ND2	1.89	0.60
1:A:24:PRO:CD	1:A:205:GLU:H	2.14	0.60
1:B:24:PRO:HA	1:B:204:ARG:CZ	2.31	0.60
1:A:152:ALA:HA	1:A:172:VAL:HG21	1.84	0.59
1:A:162:ARG:HH11	1:B:210:ASN:ND2	1.89	0.59
1:A:93:SER:HB3	2:A:235:HOH:O	2.03	0.59
1:A:24:PRO:HD3	1:A:205:GLU:N	2.17	0.59
1:A:68:VAL:HG23	1:B:73:PRO:HG3	1.84	0.59
1:A:24:PRO:CD	2:A:229:HOH:O	2.35	0.58
1:A:155:GLN:NE2	1:A:155:GLN:N	2.52	0.58
1:B:163:ALA:O	1:B:166:ASP:HB2	2.03	0.57
1:A:144:ARG:NH2	1:A:179:GLN:O	2.33	0.57
1:A:54:ARG:HD2	2:A:244:HOH:O	2.03	0.57
1:B:148:LEU:HD23	1:B:172:VAL:HG12	1.86	0.57
1:B:111:THR:OG1	1:B:122:HIS:HD2	1.88	0.56
1:A:168:GLN:HE22	1:B:223:GLN:HE22	1.53	0.56
1:A:179:GLN:HE21	1:A:179:GLN:HA	1.69	0.56
1:A:68:VAL:CG2	1:B:73:PRO:HD3	2.36	0.56
1:B:144:ARG:NH2	1:B:179:GLN:O	2.39	0.56
1:A:144:ARG:NH1	2:A:245:HOH:O	2.39	0.56
1:B:210:ASN:HA	2:B:234:HOH:O	2.06	0.55
1:A:77:SER:H	1:B:120:GLN:HE22	1.56	0.54
1:B:155:GLN:CG	1:B:156:SER:H	2.21	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:GLN:H	1:B:223:GLN:HE22	1.54	0.53
1:A:111:THR:OG1	1:A:122:HIS:HD2	1.91	0.53
1:A:120:GLN:NE2	1:B:77:SER:OG	2.42	0.52
1:B:224:PRO:C	1:B:226:SER:H	2.12	0.52
1:B:32:LEU:HG	1:B:212:ILE:HD12	1.92	0.52
1:B:211:ARG:HH12	1:B:226:SER:CA	2.20	0.52
1:A:162:ARG:NH1	1:B:210:ASN:ND2	2.51	0.51
1:A:155:GLN:H	1:A:155:GLN:CD	2.13	0.51
1:A:152:ALA:HB3	1:B:208:MET:CE	2.42	0.50
1:A:164:GLN:O	1:A:168:GLN:HG3	2.11	0.50
1:A:163:ALA:O	1:A:166:ASP:HB2	2.12	0.50
1:A:179:GLN:NE2	1:A:180:ASP:N	2.60	0.49
1:A:171:GLU:HA	1:A:174:ARG:NH1	2.28	0.48
1:B:49:LEU:HD12	1:B:114:TRP:HZ2	1.78	0.48
1:A:24:PRO:HA	1:A:204:ARG:CZ	2.43	0.48
1:B:83:LEU:HB3	2:B:257:HOH:O	2.12	0.48
1:A:70:ASP:HB2	2:A:266:HOH:O	2.11	0.48
1:A:144:ARG:O	1:A:145:GLY:C	2.52	0.47
1:B:204:ARG:H	1:B:209:HIS:CD2	2.27	0.47
1:A:211:ARG:HH12	1:A:226:SER:CA	2.13	0.47
1:B:25:GLU:OE2	1:B:25:GLU:HA	2.14	0.47
1:B:27:ASP:OD2	1:B:29:CME:HB3	2.14	0.47
1:B:46:ARG:HG3	1:B:83:LEU:HD13	1.96	0.47
1:A:154:GLN:N	1:B:223:GLN:NE2	2.62	0.47
1:A:179:GLN:NE2	1:A:180:ASP:H	2.12	0.46
1:B:61:ASN:HD22	1:B:61:ASN:N	2.05	0.46
1:A:154:GLN:HB3	1:A:157:ARG:HB3	1.98	0.46
1:A:152:ALA:HB3	1:B:208:MET:HE1	1.98	0.46
1:B:76:ARG:N	1:B:76:ARG:HD2	2.30	0.46
1:A:210:ASN:HD21	1:B:162:ARG:HH11	1.62	0.46
1:B:24:PRO:CD	1:B:205:GLU:HB2	2.42	0.46
1:B:210:ASN:HD22	1:B:221:ARG:HH12	1.64	0.45
1:A:207:ARG:HB2	1:B:149:GLY:HA3	1.97	0.45
1:B:32:LEU:CG	1:B:212:ILE:HD12	2.46	0.45
1:A:149:GLY:HA2	1:B:208:MET:HE3	1.99	0.45
1:B:111:THR:OG1	1:B:122:HIS:CD2	2.68	0.45
1:B:156:SER:O	1:B:157:ARG:C	2.54	0.45
1:A:76:ARG:HD2	1:A:76:ARG:N	2.32	0.44
1:A:24:PRO:CG	2:A:229:HOH:O	2.62	0.44
1:B:24:PRO:HD3	1:B:205:GLU:H	1.83	0.44
1:B:148:LEU:HD23	1:B:172:VAL:CG1	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:PRO:HD3	2:B:228:HOH:O	2.16	0.44
1:B:131:SER:O	1:B:135:ILE:HG13	2.18	0.44
1:A:120:GLN:HB2	1:A:201:TRP:HB3	2.00	0.43
1:A:155:GLN:N	1:A:155:GLN:CD	2.69	0.43
1:B:162:ARG:NH2	1:B:166:ASP:OD2	2.34	0.43
1:B:210:ASN:HD22	1:B:221:ARG:NH1	2.17	0.43
1:A:139:TRP:O	1:A:142:ARG:HB2	2.19	0.43
1:A:54:ARG:CD	2:A:244:HOH:O	2.66	0.42
1:A:77:SER:HB2	1:B:120:GLN:NE2	2.34	0.42
1:A:207:ARG:HD3	1:A:207:ARG:HA	1.67	0.42
1:A:179:GLN:NE2	1:A:179:GLN:CA	2.83	0.42
1:B:90:PHE:CZ	1:B:191:TYR:HB2	2.54	0.42
1:B:127:VAL:HA	1:B:192:ARG:O	2.19	0.42
1:B:124:GLN:HG2	1:B:197:ILE:HB	2.02	0.42
1:B:151:TRP:CD1	1:B:185:PRO:HG3	2.55	0.42
1:B:120:GLN:HB2	1:B:201:TRP:HB3	2.01	0.41
1:A:144:ARG:HA	1:A:147:GLN:HB2	2.01	0.41
1:B:210:ASN:ND2	1:B:221:ARG:HH12	2.18	0.41
1:B:130:VAL:HG23	1:B:135:ILE:HG13	2.01	0.41
1:A:68:VAL:HG21	1:B:73:PRO:HD3	2.02	0.41
1:A:82:ILE:HG12	1:A:138:TYR:CZ	2.56	0.41
1:B:24:PRO:CD	1:B:205:GLU:H	2.34	0.41
1:A:24:PRO:HD2	1:A:205:GLU:HB2	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ASP:OD2	1:B:54:ARG:NH2[4_454]	2.16	0.04

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	200/246 (81%)	187 (94%)	9 (4%)	4 (2%)	7 12
1	B	200/246 (81%)	191 (96%)	5 (2%)	4 (2%)	7 12
All	All	400/492 (81%)	378 (94%)	14 (4%)	8 (2%)	7 12

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	155	GLN
1	B	157	ARG
1	A	156	SER
1	B	155	GLN
1	B	217	GLY
1	B	25	GLU
1	A	217	GLY
1	A	113	PRO

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	164/200 (82%)	146 (89%)	18 (11%)	6 12
1	B	164/200 (82%)	150 (92%)	14 (8%)	10 21
All	All	328/400 (82%)	296 (90%)	32 (10%)	8 15

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	LEU
1	A	53	GLN
1	A	54	ARG
1	A	58	SER
1	A	61	ASN
1	A	76	ARG
1	A	83	LEU
1	A	98	LYS

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Mol	Chain	Res	Type
1	A	117	LEU
1	A	144	ARG
1	A	154	GLN
1	A	155	GLN
1	A	156	SER
1	A	157	ARG
1	A	162	ARG
1	A	179	GLN
1	A	181	GLN
1	A	207	ARG
1	B	32	LEU
1	B	42	LEU
1	B	61	ASN
1	B	76	ARG
1	B	83	LEU
1	B	98	LYS
1	B	117	LEU
1	B	132	THR
1	B	159	VAL
1	B	162	ARG
1	B	175	ARG
1	B	179	GLN
1	B	216	ASN
1	B	226	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	ASN
1	A	120	GLN
1	A	122	HIS
1	A	124	GLN
1	A	154	GLN
1	A	155	GLN
1	A	179	GLN
1	A	209	HIS
1	A	210	ASN
1	A	223	GLN
1	B	61	ASN
1	B	120	GLN
1	B	122	HIS
1	B	154	GLN

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Mol	Chain	Res	Type
1	B	167	ASN
1	B	179	GLN
1	B	209	HIS
1	B	210	ASN
1	B	223	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	CME	A	29	1	8,9,10	0.90	0	5,9,11	1.89	1 (20%)
1	CME	B	29	1	8,9,10	0.95	1 (12%)	5,9,11	1.92	1 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CME	A	29	1	-	1/5/8/10	-
1	CME	B	29	1	-	1/5/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	29	CME	CB-SG	-2.14	1.74	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	29	CME	CB-SG-SD	3.57	113.08	103.82
1	A	29	CME	CB-SG-SD	3.22	112.16	103.82

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	29	CME	CZ-CE-SD-SG
1	B	29	CME	SD-CE-CZ-OH

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	29	CME	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	30:GLY	C	31:ASP	N	1.19
1	A	174:ARG	C	175:ARG	N	1.19
1	B	28:GLY	C	29:CME	N	1.18

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	202/246 (82%)	0.08	7 (3%)	44	47	27, 35, 46, 60	0
1	B	202/246 (82%)	-0.06	5 (2%)	57	61	26, 34, 47, 61	0
All	All	404/492 (82%)	0.01	12 (2%)	50	53	26, 35, 47, 61	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	24	PRO	3.5
1	A	180	ASP	3.5
1	A	177	ALA	3.5
1	B	180	ASP	3.2
1	A	216	ASN	2.9
1	B	217	GLY	2.8
1	B	156	SER	2.8
1	B	24	PRO	2.7
1	A	54	ARG	2.7
1	A	174	ARG	2.6
1	B	216	ASN	2.6
1	A	181	GLN	2.5

6.2 Non-standard residues in protein, DNA, RNA chains (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
1	CME	A	29	10/11	0.89	0.20	30,34,57,61	0
1	CME	B	29	10/11	0.91	0.17	52,55,66,66	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.